

# Mathematical Modeling and Analysis

## M-adaptation for acoustic wave equation

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Numerical modeling of wave propagation is essential for a large number of applied problems in acoustics, elasticity, and electromagnetics. The acoustic equation is one of the simplest examples of equation modeling wave propagation. For long integration times, the dominant contributions to an error in the solution come from such numerical artifacts as numerical dispersion and numerical anisotropy. The *numerical dispersion* is the phenomenon in which the propagation velocity of the wave in the numerical scheme depends on its wavelength, while in the continuum problem there is no such dependence. Typically, the effect of the numerical dispersion is greater on under-resolved waves with ten or less points per wavelength, making them travel slower than in the physical problem. As a consequence, the wave does not just arrive at a wrong time (which could be compensated by time re-scaling) but it also has a highly distorted profile. The *numerical anisotropy* is the dependence of the numerical velocity of the wave on its orientation with respect to the mesh. For a 2D acoustic wave equation we developed an adaptation technique, dubbed m-adaptation, that selects an optimal member of a rich parameterized family of second order methods with smallest (fourth order) dispersion and (sixth order) anisotropy.

The semi-discrete form of the acoustic wave equation in the time domain formulation is

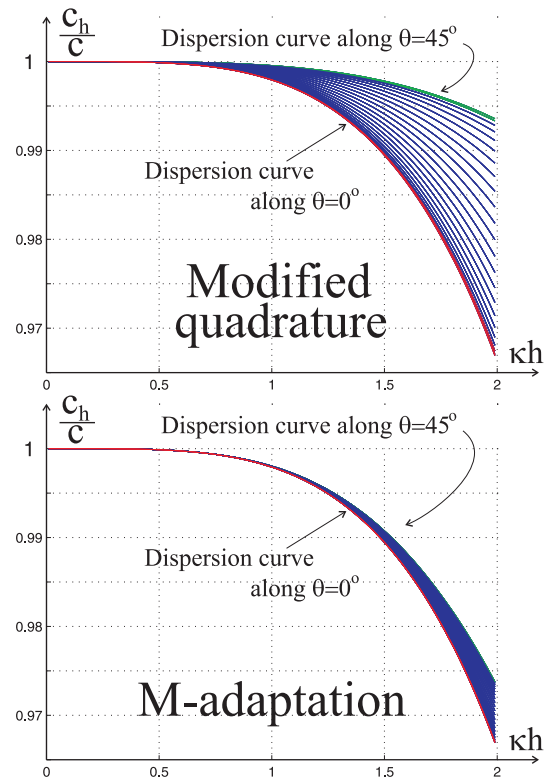
$$Mu_{tt} = Au, \quad (1)$$

where the mass and stiffness matrices  $M$  and  $A$  are assembled from elemental matrices  $M_E$  and  $A_E$ .

Since the mass matrix  $M$  has to be inverted on every time step, the explicit time discretization of equation (1) is computationally efficient only when the inverse  $M^{-1}$  is easy to compute. One of the approaches is to replace the mass

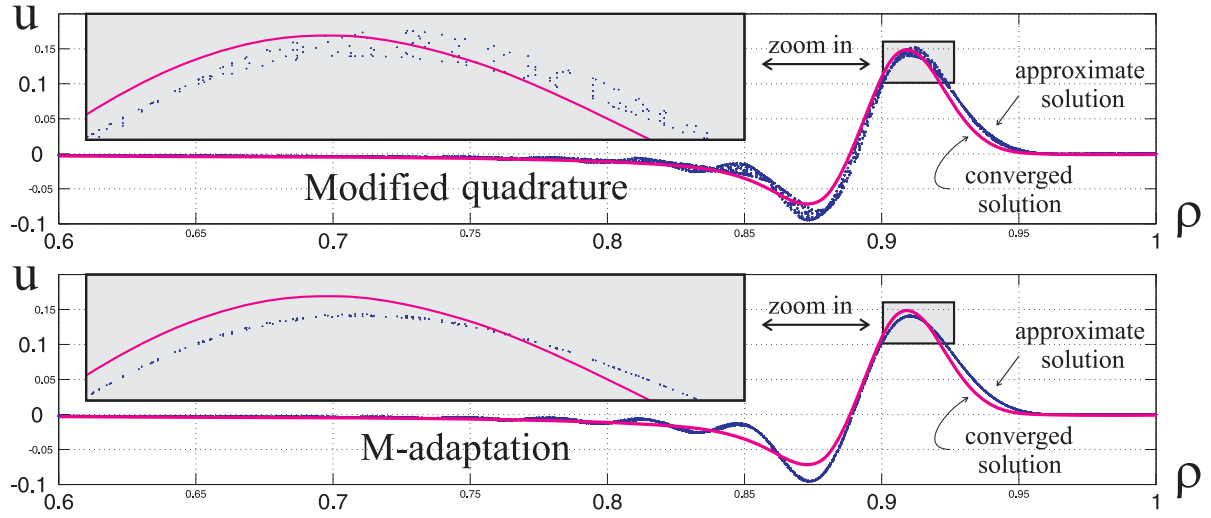
matrix  $M$  with a diagonal matrix  $D$  by lumping non-diagonal entries to the diagonal. This does not change the order of the numerical scheme but may lead to undesirable increase of numerical dispersion. Another approach (see e.g. [2]) is to replace the inverse  $M^{-1}$  with the product  $D^{-1}MD^{-1}$ , where the inverse is taken only for the diagonal matrix  $D$ . Similar to lumping, this approach does not change the order of the numerical scheme but may also result in the increase of the numerical dispersion. To compensate for the possible increase of the dispersion one can modify the stiffness and the mass matrices  $A$  and  $M$  using modified quadrature rules as it is done in [2].

In the m-adaptation approach, we consider a parameterized Mimetic Finite Difference (MFD)



A more narrow band of values in the dispersion curves for the m-adaptation method (bottom) compared to the modified quadrature method [2] (top) for various angles  $\theta$  between the planar wave and the mesh axis for the Courant number  $\frac{c\Delta t}{h} = 0.75$  indicates smaller anisotropy.

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Displacement as a function of the distance from the origin at time  $T = 0.9$  obtained using the modified quadrature method (top) and the m-adaptation method (bottom) for a Gaussian initial displacement data.

family of numerical schemes from which we select a member with the smallest numerical dispersion and anisotropy (see [1] for details). The parameters in the MFD family appear through the elemental mass and stiffness matrices  $M_E^{\text{MFD}}$  and  $A_E^{\text{MFD}}$ , respectively. The elemental mass matrix  $M_E^{\text{MFD}}$  on a square element  $E$  depends on two parameters  $m_1, m_2$  while the elemental stiffness matrix  $A_E$  depends on one parameter  $\zeta$ .

The MFD family parameterized by  $(m_1, m_2, \zeta)$  contains a large number of known methods as special cases, e.g.: standard Finite Difference (FD), rotated FD, weighted combination of standard and rotated FD, Finite Element (FE) with lumped mass matrix, and modified quadrature method of Guddati and Yue [2]. Moreover, compared with the later method, the MFD family is richer – it contains one extra parameter.

For the acoustic wave equation in 2D the optimal parameters  $(m_1, m_2, \zeta)$  can be selected based on the von-Neumann analysis. One obtains a local dispersion equation relating the numerical velocity of the wave  $c_h$  with its wave number  $\kappa$ , mesh size  $h$ , and the parameters  $(m_1, m_2, \zeta)$ . Expanding the error between the physical and numerical velocities of the wave,  $c - c_h$ , in powers of wave resolution number,  $\kappa h$ , we select the

parameters  $(m_1, m_2, \zeta)$  to eliminate the error at the leading powers of  $\kappa h$ . As a result of m-adaptation the numerical velocity  $c_h$  is accurate to the fourth order in dispersion (as in [2]) and to the sixth order in the anisotropy (versus fourth order in [2]).

In the future we plan to develop the m-adaptation technique for higher order schemes on general meshes and for elastic wave equations. The potential of m-adaptation is high as with increase of the order of the scheme and/or the number of vertices in the element the number of free parameters grows quadratically. This may lead to a dramatic improvement in the dispersion and anisotropy of the optimal scheme.

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## References

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